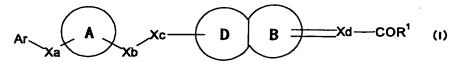
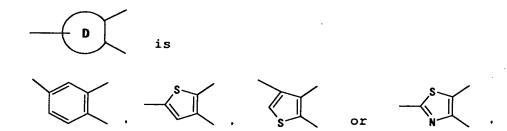
## Claims

1. A compound represented by the formula:



- wherein Ar is an optionally substituted cyclic group,
  ring A is a ring optionally further substituted (provided that
  the ring is not thiazole, oxazole, imidazole and pyrazole),
  Xa and Xb are each independently a bond or a spacer having a
  main chain of 1 to 5 atom(s),
- 10 Xc is O, S, SO or  $SO_2$ ,



ring B is a 5- to 7-membered ring,

Xd is a bond, CH or  $CH_2$ ,

 $\cdots$  is a single bond when Xd is a bond or  $CH_2$ , or a double bond when Xd is  $CH_2$ .

 ${\ensuremath{\mathsf{R}}}^1$  is an optionally substituted hydroxy group, provided that

- (i) when ring A is benzene, the cyclic group represented by Ar is not a quinolinyl group,
- 20 (ii) when ring B is a 5- to 7-membered aromatic ring, the ring represented by ring A is not thiophene and furan,
  - (iii) when ring B is benzene, the ring represented by ring A is not 5-membered aromatic heterocycle, and
  - (iv) when ring B is cyclohexane, Xd is not a bond,
- 25 provided that

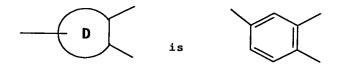
[6-(4-biphenylyl)methoxy-2-tetralin]acetic acid;
methyl [6-(4-biphenylyl)methoxy-2-tetralin]acetate;

[7-(4-biphenyly1)methoxy-1,2,3,4-tetrahydro-2-oxo-3-quinoline]acetic acid; and methyl [7-(4-biphenyly1)methoxy-1,2,3,4-tetrahydro-2-oxo-3-quinoline]acetate are excluded,

- 5 or a salt thereof.
  - 2. A prodrug of the compound of claim 1.
- 3. The compound of claim 1, wherein the cyclic group represented by Ar is an aromatic hydrocarbon group.
  - 4. The compound of claim 1, wherein Xa is a bond.
  - 5. The compound of claim 1, wherein ring A is benzene.

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- 6. The compound of claim 1, wherein Xb is  $-CH_2-$ .
- 7. The compound of claim 1, wherein Xc is 0.
- 20 8. The compound of claim 1, wherein



9. The compound of claim 1, wherein ring B is a 5- to 7-membered non-aromatic ring.

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- 10. The compound of claim 9, wherein ring B is cyclopentane or tetrahydrofuran.
- 11. The compound of claim 1, wherein Xd is  $CH_2$ .

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12. The compound of claim 1, wherein R1 is a hydroxy group.

13. The compound of claim 1, which is represented by the formula:

$$Ar^1 - Xa^1 - A^2$$
 $CH_2COOH$ 
 $B^2$ 

wherein  $Ar^1$  is an optionally substituted phenyl group or optionally substituted indanyl group,  $Xa^1$  is a bond or a spacer having a main chain of 1 to 5 atom(s), ring  $A^2$  is a benzene ring optionally further substituted, and ring  $B^2$  is a 5- to 7-membered ring.

10 14. The compound of claim 1, which is represented by the formula:

$$Ar^2 - Xa^2 - A^3$$

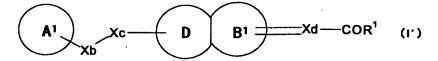
$$CH_2COOH$$

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wherein  $Ar^2$  is an optionally substituted thiazolyl group,  $Xa^2$  is a bond or a spacer having a main chain of 1 to 5 atom(s), ring  $A^3$  is a benzene ring optionally further substituted, and ring  $B^2$  is a 5- to 7-membered ring.

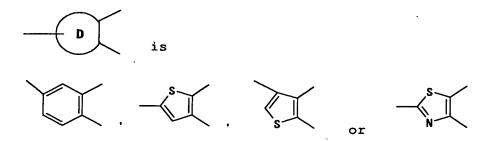
- 15. A pharmaceutical agent comprising the compound of claim 1 or a prodrug thereof.
- 16. The pharmaceutical agent of claim 15, which is an agent for the prophylaxis or treatment of diabetes.
- 17. An insulin secretagogue comprising the compound of claim 1 25 or a prodrug thereof.

18. A GPR40 receptor function modulator comprising a compound represented by the formula:



wherein ring A<sup>1</sup> is an optionally substituted ring,

5 Xb is a bond or a spacer having a main chain of 1 to 5 atom(s), Xc is O, S, SO or  $SO_2$ ,



ring  $B^1$  is a 5- to 7-membered non-aromatic ring, Xd is a bond, CH or  $CH_2$ ,

- 10  $\cdots$  is a single bond when Xd is a bond or  $CH_2$ , or a double bond when Xd is CH, and  $R^1$  is an optionally substituted hydroxy group, or a salt thereof, or a prodrug thereof.
- 15 19. A method of modulating a GPR40 receptor function in a mammal, which comprises administering an effective amount of the compound of claim 18 or a prodrug thereof to the mammal.
- 20. Use of the compound of claim 18 or a prodrug thereof for the production of a GPR40 receptor function modulator.